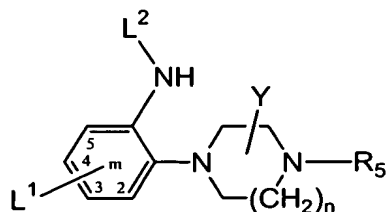


What is claimed is:

1. A compound of formula (I):



formula (I)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:

5 L¹ is a substituent moiety having a variable position “m”, wherein “m” represents a carbon atom number corresponding to a point of attachment for the L¹ substituent moiety on the anilino ring of formula (I);

L¹ is selected from the group consisting of R_{1b}, R₂-C(O), R_{1a}-SO₂ and R_{1a}-O(O)C-;

10 R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

15 R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

20 L² is selected from the group consisting of R₃-C(O)-, R₄-SO₂-, R₆-NHC(S)- and R₆-NHC(O)-;

R₃ is selected from the group consisting of

(a) C₁₋₈alkyl optionally substituted with one or more substituents independently
25 selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
wherein said aryl is optionally substituted with one or more substituents

independently selected from the group consisting of C₁₋₈alkyl,
C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano,
halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom
with C₁₋₈alkyl, and optionally and independently substituted on one or
more carbon atoms with a substituent selected from the group consisting
of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino,
di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(b) aryl optionally substituted with one or more substituents independently selected
from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
and,

(c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and
optionally and independently substituted on one or more carbon atoms with a
substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and
aryl, wherein said aryl is optionally substituted with one or more substituents
independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy,
amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and
nitro;

R₄ is selected from the group consisting of

(d) C₁₋₈alkyl optionally substituted with one or more substituents independently
selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino,
di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,

(e) aryl optionally substituted with one or more substituents independently selected
from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or more substituents independently selected
from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and, wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (I); and, n is an integer from 1 to 2.

2. The compound of claim 1, wherein when L^2 is $R_3-C(O)-$ and R_3 is selected from the group consisting of unsubstituted C_{1-8} alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then L^1 is $R_2-C(O)$.
- 5 3. The compound of claim 1, wherein when L^2 is $R_3-C(O)-$ and R_3 is selected from the group consisting of unsubstituted C_{1-8} alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then R_5 is C_{1-8} alkyl optionally substituted with one or more optionally substituted aryl substituents.
- 10 4. The compound of claim 1, wherein when L^2 is R_4-SO_2- and R_4 is unsubstituted C_{1-8} alkyl, then L^1 is $R_2-C(O)$, wherein R_2 is substituted or unsubstituted heterocyclyl.
- 15 5. The compound of claim 1, wherein when L^2 is R_4-SO_2- and R_4 is unsubstituted C_{1-8} alkyl, then R_5 is C_{1-8} alkyl optionally substituted with one or more optionally substituted aryl substituents.
- 20 6. The compound of claim 1, wherein when L^1 is selected from the group consisting of R_{1b} and $R_{1a}-O(O)C-$, then L^2 is $R_6-NHC(O)-$, wherein R_6 is substituted or unsubstituted aryl.
- 25 7. The compound of claim 1, wherein when L^1 is selected from the group consisting of R_{1b} and $R_{1a}-O(O)C-$, then R_5 is C_{1-8} alkyl optionally substituted with one or more optionally substituted aryl substituents.
- 30 8. The compound of claim 1, wherein R_{1a} is C_{1-8} alkyl optionally substituted with one or two substituents independently selected from the group consisting of C_{1-8} alkoxy, amino, mono(C_{1-8})alkylamino, di(C_{1-8})alkylamino, halogen and hydroxy;
- R_{1b} is C_{1-8} alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C_{1-8})alkylamino, di(C_{1-8})alkylamino, halogen and hydroxy;

R₃ is selected from the group consisting of

- 5 (a) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
10 wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 15 (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- 20 (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy,
25 amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- 30 (d) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,

mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,

5 mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

(f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently
10 selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;

wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl,
15 C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting
20 of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
25 and,

(h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
and,

30

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro,

C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

9. The compound of claim 1, wherein when L² is R₃-C(O)- and R₃ is selected from the group consisting of unsubstituted C₁₋₈alkyl, substituted aryl, unsubstituted aryl, substituted heteroaryl and unsubstituted heteroaryl, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
10. The compound of claim 1, wherein when L² is R₄-SO₂- and R₄ is unsubstituted C₁₋₈alkyl, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
11. The compound of claim 1, wherein when L¹ is selected from the group consisting of R_{1b} and R_{1a}-O(O)C-, then R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
12. The compound of claim 1, wherein
R_{1a} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy;
- R_{1b} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy;
- R₂ is piperazinyl optionally substituted on a nitrogen atom with C₁₋₄alkyl;
- L² is selected from the group consisting of R₃-C(O)-, R₄-SO₂- and R₆-NHC(O)-;
- R₃ is selected from the group consisting of
- (a) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;

- (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,
- (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano,

halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

(h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is absent;

m is an integer from 3 to 4 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (I); and, n is 1.

13. The compound of claim 12, wherein R_{1a} is C₁₋₄alkyl;

R_{1b} is hydroxy(C₁₋₄)alkyl-;

R₃ is selected from the group consisting of

(a) C₁₋₄alkyl;

(b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,

(c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and aryl,

wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

5

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl; and,
- (e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, 10 mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

15

R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

20

R₅ is selected from the group consisting of

- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- (g) C₃₋₈cycloalkyl; and,
- (h) aryl.

25

14. The compound of claim 13, wherein R₃ is selected from the group consisting of

- (a) C₁₋₄alkyl;
- (b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen; and,
- (c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl and phenyl; 30 wherein said phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen;

R₄ is selected from the group consisting of

- (d) C₁₋₄alkyl; and,
- (e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen;

5

R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, amino, halogen and hydroxy; and,

10 R₅ is C₁₋₄alkyl optionally substituted with one or two phenyl substituents, wherein phenyl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, amino, halogen and hydroxy.

15 15. The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.

20 16. The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl optionally substituted with one substituent selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.

17. The compound of claim 1, wherein R_{1a} is C₁₋₄alkyl.

25 18. The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino and hydroxy.

30 19. The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally substituted with one substituent selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy.

20. The compound of claim 1, wherein R_{1b} is C₁₋₄alkyl optionally substituted with hydroxy.

21. The compound of claim 1, wherein R₂ is piperazinyl optionally substituted on a nitrogen atom with C₁₋₄alkyl.
- 5 22. The compound of claim 1, wherein L² is R₃-C(O)-.
23. The compound of claim 22, wherein R₃ is selected from the group consisting of
- 10 (a) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;
- (b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy;
- 15 (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₄alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
- 20 24. The compound of claim 22, wherein R₃ is selected from the group consisting of
- (a) C₁₋₄alkyl;
- 25 (b) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, halogen and hydroxy; and,
- (c) furyl optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and
- 30

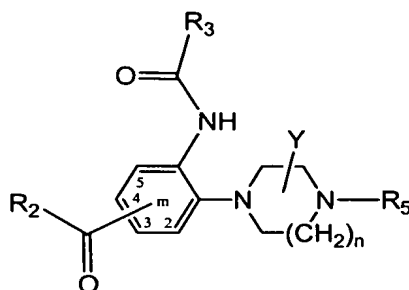
nitro.

25. The compound of claim 22, wherein R_3 is selected from the group consisting of
- (a) C_{1-4} alkyl;
- 5 (b) phenyl optionally substituted with one or two substituents independently
selected from the group consisting of C_{1-4} alkyl and halogen; and,
- (c) furyl optionally and independently substituted on one or two carbon atoms with
a substituent selected from the group consisting of C_{1-4} alkyl and phenyl;
wherein said phenyl is optionally substituted with one or two substituents
10 independently selected from the group consisting of C_{1-4} alkyl and halogen.
26. The compound of claim 1, wherein L^2 is $R_3-C(O)-$ and R_5 is C_{1-8} alkyl optionally
substituted with one or two optionally substituted aryl substituents.
- 15 27. The compound of claim 26, wherein R_3 is selected from the group consisting of
- (a) C_{1-4} alkyl optionally substituted with one or two substituents independently
selected from the group consisting of C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino,
di(C_{1-4})alkylamino, hydroxy, aryl and heteroaryl;
- (b) aryl optionally substituted with one or two substituents independently selected
20 from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino,
mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, halogen and hydroxy;
- (c) heteroaryl optionally substituted on a secondary amine atom with C_{1-4} alkyl, and
optionally and independently substituted on one or two carbon atoms with a
substituent selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy, amino,
25 mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy, nitro and
aryl, wherein said aryl is optionally substituted with one or two substituents
independently selected from the group consisting of C_{1-4} alkyl, C_{1-4} alkoxy,
amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and
nitro.
30
28. The compound of claim 1, wherein L^2 is R_4-SO_2- .
29. The compound of claim 28, wherein R_4 is selected from the group consisting of

- (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
30. The compound of claim 28, wherein R₄ is selected from the group consisting of
- (d) C₁₋₄alkyl; and,
- (e) phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl and halogen.
31. The compound of claim 1, wherein L² is R₄-SO₂- and R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.
32. The compound of claim 31, wherein R₄ is selected from the group consisting of
- (d) C₁₋₄alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
- (e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
33. The compound of claim 1, wherein L² is R₆-NHC(O)-.
34. The compound of claim 33, wherein R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
35. The compound of claim 1, wherein L² is R₆-NHC(O)- and R₅ is C₁₋₈alkyl optionally substituted with one or two optionally substituted aryl substituents.

36. The compound of claim 35, wherein R₆ is phenyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
- 5
37. The compound of claim 1, wherein R₅ is selected from the group consisting of
- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, 10 mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; 15 and,
- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro.
- 20 38. The compound of claim 1, wherein R₅ is selected from the group consisting of
- (f) C₁₋₄alkyl optionally substituted with one or two aryl substituents, wherein aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₄alkyl, C₁₋₄alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 25 (g) C₃₋₈cycloalkyl; and,
- (h) aryl.
39. The compound of claim 1, wherein R₅ is C₁₋₄alkyl optionally substituted with one or two phenyl substituents, wherein phenyl is optionally substituted with 30 one or two substituents independently selected from the group consisting of C₁₋₄alkyl, amino, halogen and hydroxy.

40. The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ia):



formula (Ia)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:
R₂-C(O)- is a substituent moiety having a variable position “m”, wherein “m”

- 5 represents a carbon atom number corresponding to a point of attachment for the
R₂-C(O)- substituent moiety on the anilino ring of formula (Ia);

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

- 10 R₃ is selected from the group consisting of

(a) C₁₋₈alkyl optionally substituted with one or more substituents independently
selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino,
di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;

wherein said aryl is optionally substituted with one or more substituents

- 15 independently selected from the group consisting of C₁₋₈alkyl,
C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano,
halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom

- 20 with C₁₋₈alkyl, and optionally and independently substituted on one or
more carbon atoms with a substituent selected from the group consisting
of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino,
di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(b) aryl optionally substituted with one or more substituents independently selected
from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
25 mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
and,

- (c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 10 R₅ is selected from the group consisting of
- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- 15 wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- 20 wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 25 (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 30

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ia); and, n is an integer from 1 to 2.

41. The compound of claim 40, wherein R₃ is selected from the group consisting of
(a) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl;

wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(b) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

(c) heteroaryl optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro and aryl, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy,

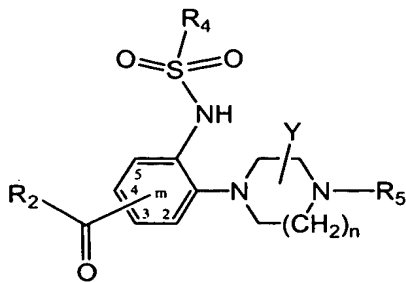
amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- 5 (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- 10 wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- 15 wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 20 (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- 25 (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, 30 mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

42. The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ib):



formula (Ib)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:
R₂-C(O)- is a substituent moiety having a variable position “m”, wherein “m”

- 5 represents a carbon atom number corresponding to a point of attachment for the
R₂-C(O)- substituent moiety on the anilino ring of formula (Ib);

R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

- 10 R₄ is selected from the group consisting of

- (d) C₁₋₈alkyl optionally substituted with one or more substituents independently
selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino,
di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,
(e) aryl optionally substituted with one or more substituents independently selected
15 from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said
20 aryl is optionally substituted with one or more substituents independently
selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl'
and heteroaryl;
wherein said aryl' is optionally substituted with one or more substituents
25 independently selected from the group consisting of C₁₋₈alkyl,
C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano,

halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

(h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the R₂-C(O)- substituent moiety on the anilino ring of formula (Ib); and, n is an integer from 1 to 2.

43. The compound of claim 42, wherein R₄ is selected from the group consisting of

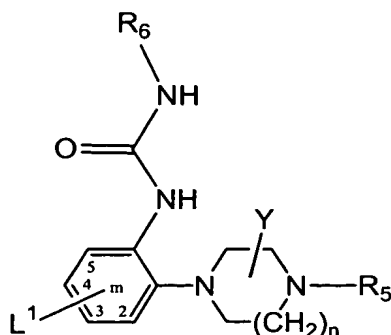
(d) C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, hydroxy, aryl and heteroaryl; and,

(e) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

- (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl; wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and, wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- (g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

44. The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Ic):



formula (Ic)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:

- L^1 is a substituent moiety having a variable position “m”, wherein “m” represents a
5 carbon atom number corresponding to a point of attachment for the L^1
substituent moiety on the anilino ring of formula (Ic);

L^1 is selected from the group consisting of R_{1b} , R_{1a} -SO₂- and R_{1a} -O(O)C-;

- 10 R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently
selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino,
di(C₁₋₈)alkylamino, halogen and hydroxy;

- 15 R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently
selected from the group consisting of amino, mono(C₁₋₈)alkylamino,
di(C₁₋₈)alkylamino, halogen and hydroxy;

- 20 R_6 is aryl optionally substituted with one or more substituents independently selected
from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,
mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

- R_5 is selected from the group consisting of
(f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said
aryl is optionally substituted with one or more substituents independently
25 selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,

mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;

wherein said aryl' is optionally substituted with one or more substituents

independently selected from the group consisting of C₁₋₈alkyl,

5 C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

wherein said heteroaryl is optionally substituted on a secondary amine atom

with C₁₋₈alkyl, and optionally and independently substituted on one or

more carbon atoms with a substituent selected from the group consisting

10 of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or more substituents

independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy,

amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and

15 nitro; and,

(h) aryl optionally substituted with one or more substituents independently selected

from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino,

mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

20 Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;

25

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (Ic); and, n is an integer from 1 to 2.

30 45. The compound of claim 44, wherein

R_{1a} is C₁₋₈alkyl optionally substituted with one or two substituents independently

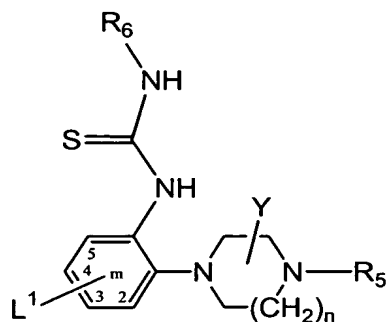
selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino,

di(C₁₋₈)alkylamino, halogen and hydroxy;

- R_{1b} is C_{1-8} alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C_{1-8})alkylamino, di(C_{1-8})alkylamino, halogen and hydroxy;
- 5
- R_6 is aryl optionally substituted with one or two substituents independently selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro;
- 10
- R_5 is selected from the group consisting of
- (f) C_{1-8} alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- 15
- wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro; and,
- 20
- wherein said heteroaryl is optionally substituted on a secondary amine atom with C_{1-8} alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro;
- 25
- (g) C_{3-8} cycloalkyl optionally substituted with one or two substituents independently selected from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or two substituents independently selected
- 30
- from the group consisting of C_{1-8} alkyl, C_{1-8} alkoxy, amino, mono(C_{1-4})alkylamino, di(C_{1-4})alkylamino, cyano, halogen, hydroxy and nitro; and,

Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

46. The compound of claim 1, wherein the compound of formula (I) is selected from a compound of formula (Id):



formula (Id)

and enantiomers, diastereomers and pharmaceutically acceptable salts thereof, wherein:

10 L¹ is a substituent moiety having a variable position “m”, wherein “m” represents a carbon atom number corresponding to a point of attachment for the L¹ substituent moiety on the anilino ring of formula (Id);

15 L¹ is selected from the group consisting of R_{1b}, R₂-C(O)-, R_{1a}-SO₂- and R_{1a}-O(O)C-;

R_{1a} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

20 R_{1b} is C₁₋₈alkyl optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

25 R₂ is heterocyclyl optionally substituted on a nitrogen atom with C₁₋₈alkyl;

- R₆ is aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 5 R₅ is selected from the group consisting of
- (f) C₁₋₈alkyl optionally substituted with one or more aryl substituents, wherein said aryl is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;
- 10 wherein said aryl' is optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- 15 wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or more carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 20 (g) C₃₋₈cycloalkyl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,
- (h) aryl optionally substituted with one or more substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;
- 25
- Y is one or more optionally present C₁₋₈alkyl substituents optionally substituted with one or more substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted;
- 30

m is an integer from 2 to 5 which represents the carbon atom number corresponding to the point of attachment for the L¹ substituent moiety on the anilino ring of formula (Id); and, n is an integer from 1 to 2.

5 47. The compound of claim 46, wherein

R_{1a} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkoxy, amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

10 R_{1b} is C₁₋₈alkyl optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₈)alkylamino, di(C₁₋₈)alkylamino, halogen and hydroxy;

15 R₆ is aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

R₅ is selected from the group consisting of

20 (f) C₁₋₈alkyl optionally substituted with one or two aryl substituents, wherein said aryl is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, aryl' and heteroaryl;

25 wherein said aryl' is optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

30 wherein said heteroaryl is optionally substituted on a secondary amine atom with C₁₋₈alkyl, and optionally and independently substituted on one or two carbon atoms with a substituent selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

(g) C₃₋₈cycloalkyl optionally substituted with one or two substituents independently

selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro; and,

- 5 (h) aryl optionally substituted with one or two substituents independently selected from the group consisting of C₁₋₈alkyl, C₁₋₈alkoxy, amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy and nitro;

10 Y is one or two optionally present C₁₋₈alkyl substituents optionally substituted with one or two substituents independently selected from the group consisting of amino, mono(C₁₋₄)alkylamino, di(C₁₋₄)alkylamino, cyano, halogen, hydroxy, nitro, C₃₋₈cycloalkyl, aryl and heteroaryl, wherein said C₃₋₈cycloalkyl, aryl and heteroaryl are optionally further substituted.

48. A compound selected from the group consisting of:

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-4-methyl-benzamide;

5-(4-chlorophenyl)-*N*-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-2-methyl-3-furancarboxamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-2-furancarboxamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-propanamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-4-methyl-benzenesulfonamide;

4-chloro-*N*-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-benzenesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-1-butanesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(1-piperazinylcarbonyl)phenyl]-methanesulfonamide;

N-[2-[4-(diphenylmethyl)-1-piperazinyl]-5-(methylsulfonyl)phenyl]-*N'*-phenyl-urea.

N-[2-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-5-(hydroxymethyl)phenyl]-*N'*-phenyl-urea; and,

4-[4-[bis(4-fluorophenyl)methyl]-1-piperazinyl]-3-[[phenylamino]carbonyl]amino]-benzoic acid methyl ester.

49. A composition comprising a pharmaceutically acceptable carrier, excipient, tableting ingredient or diluent and the compound of claim 1.
50. A method of treating or preventing a disease or condition in a subject which disease or condition is affected by phospholipase modulation, which method comprises administering to the subject in need of such treatment or prevention a therapeutically effective amount of the compound of claim 1.
51. The method of claim 50, wherein the method further comprises administering to the subject in need of such treatment or prevention a therapeutically effective amount of the composition of claim 49.
52. A method of treating or ameliorating an inflammatory disorder in a subject in need thereof comprising administering to the subject a therapeutically effective amount of the compound of claim 1.
53. The method of claim 52, wherein the method further comprises administering to the subject a therapeutically effective amount of the composition of claim 49.
54. A method of treating or ameliorating restenosis in a subject in need thereof comprising administering to the subject a therapeutically effective amount of the compound of claim 1 by impregnating the therapeutically effective amount of said compound on the surface of a medical device and administering the medical device to the subject.
55. The method of claim 54, wherein the method further comprises a therapeutically effective amount of the composition of claim 49 impregnated on the surface of said medical device.